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Amendment to the Claims:

This listing of claims will replace all prior versions and listings of claims in the application:

Listing of Claims:

1. (Currently amended): A compound according to Formula (1) (I):

wherein:

X is selected from the group consisting of CH₂,NR₂,O,NR₂CO,CONR₂ and a bond; Y represents O, CH₂ or a bond:

Z represents H or F;

X is selected from the group consisting of: (CH₂)_n, NR2, O, NR2CO, CONR2 and a bond;

Y is selected from the group consisting of: O, CH₂ and a covalent bond;

Z is selected from the group consisting of: hydrogen and fluorine;

R is selected from the group consisting of:

 C_{2-6} alkyl optionally substituted by alkoxy, halogen, or C_{1-3} alkylsulfanyl; C_{2-6} alkenyl optionally substituted by alkoxy, halogen, or C_{1-3} alkylsulfanyl; C_{2-6} alkynyl optionally substituted by alkoxy, halogen, or C_{1-3} alkylsulfanyl; $(CH_2)_n$ — C_{3-6} carbocycle optionally substituted by alkoxy, halogen, or C_{1-3} alkylsulfanyl; and $(CH_2)_n$ — R_3 ; {where R_3 is phenyl, furan, benzofuran, thiophene, benzothiophene, tetrahydrofuran, tetrahydropyran, dioxane, 1,4-benzodioxane or benzo[1,3]dioxole; R_3 is optionally substituted by one or more Cl, Br, I, C_{1-3} -alkyl optionally substituted by one to three F, or C_{1-2} -alkoxy, optionally substituted by one to three F};

R1 is selected from the group consisting of:

hydrogen, C_{1-3} substituted alkyl, C_{2-3} substituted alkenyl, C_{2-3} substituted alkynyl, $(CH_2)_n$ — C_{3-6} substituted carbocycle, aryl, heteroaryl, heterocyclic, and aminocarbonyl; provided that X is $(CH_2)_n$ when R_1 represents aminocarbonyl;

R2 represents is selected from the group consisting of:

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hydrogen, or and C₁₋₃ substituted alkyl [[,]];

R3 is selected from the group consisting of:

phenyl, furan, benzofuran, thiophene, benzothiophene, tetrahydrofuran, tetrahydropyran, dioxane, 1,4-benzodioxane and benzo[1,3]dioxole, each of which is optionally substituted by one or more substituents each independently selected from:

Cl, Br, I, C₁₋₃ alkyl optionally substituted by one to three F, and C₁₋₂ alkoxy, optionally substituted by one to three F;

X represents (CH₂)_n, NR2, O, NR2CO, CONR2 or a bond;

Y represents O, CH2 or a covalent bond;

Z represents hydrogen or fluorine; preferably fluroine; and n represents an interger integer between 0 and 2; or a pharmaceutically acceptable salt [[,]] or solvate [[,]] or physiologically functional derivative thereof.

2. (Currently amended): A compound as claimed in according to claim 1, or a pharmaceutically acceptable salt or solvate thereof, with the following absolute configuration:

wherein:

X is selected from the group consisting of CH₂,NR₂,O,NR₂CO,CONR₂ and a bond; Y represents O, CH₂ or a bond:

Z represents H or F[[;]].

- 3. (Currently amended): A compound as claimed in according to claim 1, wherein X [[=]] is CH₂ or a bond; or a pharmaceutically acceptable salt [[,]] or solvate [[,]] or physiologically functional derivative thereof.
- 4. (Currently amended): A compound as claimed in according to claim 1, wherein X [[=]] is NR2; or a pharmaceutically acceptable salt [[,]] or solvate [[,]] or physiologically functional derivative thereof.

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5. (Currently amended): A compound according to claim 3 selected from the group consisting of:

N-[(R)-2-(5-Benzo[1,3]dioxol-5-yl-[1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxyformamide;

N-[(R)-2-(5-Benzyl-[1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxyformamide;

 $N-Hydroxy-N-\{(R)-2-[5-(7-methoxy-benzofuran-2-yl)-[1,3,4]oxadiazol-2-yl]-[1,3,4]oxadiazol$

heptyl}-formamide;

N-Hydroxy-N-{(R)-2-[5-(1,2,3,4-tetrahydroquinolin-6-yl)-[1,3,4]oxadiazol-2-yl]-

heptyl}-formamide;

 $N-Hydroxy-N-\{(R)-2-[5-(1,2,3,4-tetrahydro-quinolin-8-yl)-[1,3,4]oxadiazol-2-yl]-1-(R)-2-[5-(1,2,3,4-tetrahydro-quinolin-8-yl)-[1,3,4]oxadiazol-2-yl]-1-(R)$

heptyl}-formamide;

N-Hydroxy-N-[(R)-2-(5-pyridin-3-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-

formamide;

 $N-\{(R)-2-[5-(3,4-Dichloro-benzyl)-[1,3,4] oxadiazol-2-yl]-heptyl\}-N-hydroxy-$

formamide;

 $N-\{(R)-2-[5-(3,4-Dichloro-benzyl)-[1,3,4] oxadiazol-2-yl]-heptyl\}-N-hydroxy-$

formamide:

 $N-\{(R)-2-[5-(3,4-Dichloro-benzyl)-[1,3,4] \text{ oxadiazol-} 2-yl]-heptyl\}-N-hydroxy-$

formamide:

N-Hydroxy-N-{(R)-2-[5-(2-morpholin-4-yl-ethyl)-[1,3,4]oxadiazol-2-yl]-heptyl}-

formamide;

 $N-Hydroxy-N-\{(R)-2-[5-(2-morpholin-4-yl-ethyl)-[1,3,4]oxadiazol-2-yl]-heptyl\}-$

formamide;

N-{(R)-2-[5-(2,3-Dichloro-phenoxymethyl)-[1,3,4]oxadiazol-2-yl]-heptyl}-N-

hydroxy formamide:

N-Hydroxy-N-{(R)-2-[5-(4-methoxy-phenoxymethyl)-[1,3,4]oxadiazol-2-yl]-heptyl} -

formamide;

N-((R)-2-{5-[4-(4-Acetyl-piperazin-1-yl)-phenoxymethyl]-[1,3,4]oxadiazol-2-yl}-

heptyl)-N-hydroxy-formamide;

 $N-Hydroxy-N-\{(R)-2-[5-(1-methyl-1H-pyrrol-2-ylmethyl)-[1,3,4]oxadiazol-2-yl]-$

heptyl}-formamide;

N-Hydroxy-N-[(R)-2-(5-pyridin-2-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-

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formamide;

N-Hydroxy-N-[(R)-2-(5-pyridin-4-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-

formamide;

 $N-Hydroxy-N-\{(R)-2-[5-(2,6-dichloro-benzyl)-[1,3,4]oxadiazol-2-yl]-heptyl\}$

formamide;

 $N-Hydroxy-N-\{(R)-2-[5-(1H-indol-3-ylmethyl)-[1,3,4]oxadiazol-2-yl]-heptyl\}-1-(1,3,4)oxadiazol-2-yl]-heptyl-1-(1,3,4)oxadiazo$

formamide;

N-{(R)-2-[(S)-5-(2,3-Dihydro-benzo[1,4]dioxin-2-yl)-[1,3,4]oxadiazol-2-yl]-

heptyl}-N-hydroxy-formamide;

N-[(R)-2-(5-Benzo furan-4-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxy-left (R)-2-(5-Benzo furan-4-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxy-left (R)-2-(5-Benzo furan-4-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxy-left (R)-2-(5-Benzo furan-4-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxy-left (R)-2-(5-Benzo furan-4-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxy-left (R)-2-(1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxy-left (R)-2-(1,3,4)-hydroxy-left (R)-1-(1,3,4)-hydroxy-left (

formamide;

N-Hydroxy-N-[(R)-2-(5-pyrimidin-2-yl-[1,3,4]oxadiazol-2-yl)-heptyl]-formamide;

 $N-\{(R)-2-[5-(2,3-Dihydro-benzo[d]isoxazol-3-ylmethyl)-[1,3,4]oxadiazol-2-yl]-$

heptyl}-N-hydroxy-formamide;

N-Hydroxy-N-[(R)-2-(5-phenoxymethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-formamide;

N-Hydroxy-N-{(R)-2-[(S)-5-(1,2,3,4-tetrahydro-isoquinolin-3-yl)-[1,3,4]oxadiazol-2-yl]-

heptyl}-formamide;

N-Hydroxy-N-{(R)-2-[5-(4-imidazol-1-yl-phenoxymethyl)-[1,3,4]oxadiazol-2-yl]-

heptyl}-formamide;

N-Hydroxy-N-{(R)-2-[5-(quinolin-6-yloxymethyl)-[1,3,4]oxadiazol-2-yl]-heptyl}-

formamide;

5-{(R)-1-[(Formyl-hydroxy-amino)-methyl]-hexyl}-[1,3,4]oxadiazole-2-carboxylic

acid phenylamide;

N-Hydroxy-N-[(R)-2-(5-phenylaminomethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-

formamide:

 $N-\{(R)-2-[5-(2-Chloro-benzyl)-[1,3,4] oxadiazol-2-yl]-heptyl\}-N-Hydroxy-$

formamide;

N-[(R)-2-(5-Benzyl-[1,3,4]oxadiazol-2-yl)-3-cyclohexyl-propyl]-N-hydroxyformamide.

 $N-((R)-2-\{5-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl\}-heptyl)-N-Hydroxy-level (R)-2-\{5-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl\}-heptyl)-N-Hydroxy-level (R)-2-\{5-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl\}-heptyl)-N-Hydroxy-level (R)-2-\{5-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl\}-heptyl)-N-Hydroxy-level (R)-2-\{5-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl\}-heptyl)-N-Hydroxy-level (R)-2-\{5-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl]-heptyl)-N-Hydroxy-level (R)-2-\{5-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl]-heptyl)-N-Hydroxy-level (R)-2-\{5-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl]-heptyl)-N-Hydroxy-level (R)-2-[2-(1H-Benzoimidazol-2-yl)-[2-(1H-Benzoimidazol-2-yl)-ethyl]-[1,3,4] oxadiazol-2-yl]-heptyl-1-(1H-Benzoimidazol-2-yl)-h$

formamide;

N-Hydroxy-N-{(R)-2-[5-(4-pyrimidin-2-yl-piperazin-1-ylmethyl)-[1,3,4]oxadiazol-2-yl]-

heptyl}-formamide;

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N-Hydroxy-N-{(R)-2-[5-(3-methyl-isoxazol-5-ylmethyl)-[1,3,4]oxadiazol-2-yl]-heptyl}-formamide;

N-Hydroxy-N-{(R)-2-[5-(5-methyl-2-phenyl-oxazol-4-ylmethyl)-[1,3,4]oxadiazol-2-yl]-heptyl}-formamide;

N-Hydroxy-N-(2-{5-[(4-trifluoromethyl-pyrimidin-2-ylamino)-methyl]-[1,3,4]oxadiazol-2-yl}-heptyl)-formamide;

N-{(R)-2-[5-(1H-Benzoimidazol-2-ylmethyl)-[1,3,4]oxadiazol-2-yl]-heptyl}-N-hydroxy-formamide;

N-Hydroxy-N-[(R)-2-(5-morpholin-4-ylmethyl-[1,3,4]oxadiazol-2-yl)-heptyl]-formamide; N-Hydroxy-N-((R)-2-{5-[(3-trifluoromethyl-phenylamino)-methyl]-[1,3,4]oxadiazol-2-yl}-heptyl)-formamide;

N-[(R)-2-(5-Benzyl-[1,3,4]oxadiazol-2-yl)-3-cyclopentyl-propyl]-N-hydroxyformamide; N-Hydroxy-N-[(R)-2-(5-methyl-[1,3,4]oxadiazol-2-yl)-heptyl]-formamide; and N-Hydroxy-N-((R)-2-{5-[2-(1H-indol-3-yl)-ethyl]-[1,3,4]oxadiazol-2-yl}-heptyl)-formamide; and pharmaceutically acceptable salts or solvates thereof.

6. (Currently amended): A compound according to claim 5 selected from the group consisting of:

N-Hydroxy--[(R)-2-(5-phenylamino-[1,3,4]oxadiazol-2-yl)-heptyl]-formamide;

 $N-(5-\{(R)-1-[(Formyl-hydroxy-amino)-methyl]-hexyl\}-[1,3,4]oxadiazol-2-yl)-benzamide;\\$

N-{(R)-2-[5-(Chloro-trifluoromethyl-phenylamino)-[1,3,4]oxadiazol-2-yl]-heptyl}-N-hydroxy-formamide;

N-Hydroxy-N-{(R)-2-[5-(methyl-phenyl-amino)-[1,3,4]oxadiazol-2-yl]-heptyl}-formamide; Benzo[1,3]dioxole-5-carboxylic acid (5-{(R)-1-[(formyl-hydroxy-amino)-methyl]-hexyl}-[1,3,4]oxadiazol-1,2-yl)-amide;

N-{(R)-2-[5-(3,5-Dichloro-phenylamino)-[1,3,4]oxadiazol-2-yl]-heptyl}-N-hydroxy formamide;

N-[(S)-2-Fluoro-2-(5-phenylamino-[1,3,4]oxadiazol-2-yl)-heptyl]-N-hydroxy formamide; and N-{(R)-2-[5-(2,3-Dihydro-benzo[1,4]dioxin-6-ylamino)-[1,3,4]oxadiazol-2-yl]-heptyl}-N-hydroxy-formamide; and pharmaceutically acceptable salts or solvates thereof.

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7. (Currently amended): A method of treating a bacterial infection by administering to a subject in need of such treatment a compound according to claim 1; or a pharmaceutically

acceptable salt or solvate thereof.

8. (New) A pharmaceutical composition comprising a compound according to claim 1, or a

pharmaceutically acceptable salt or solvate thereof, and one or more pharmaceutically

acceptable excipients.

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